

October 10, 2012

Mr. David Lacey Oregon Department of Environmental Quality 2020 SW Fourth Avenue, Suite 400 Portland, OR 97201-4987

**Subject:** Response to DEQ Comments

Risk Assessment, Feasibility Study, and Source Control Evaluation

Swan Island Upland Facility, Operable Unit 4

**ECSI No. 271** 

## Dear David:

This letter provides the Oregon Department of Environmental Quality (DEQ) with a response to the comments received on the Swan Island Upland Facility, Operable Unit 4 *Risk Assessment, Feasibility Study, and Source Control Evaluation* (Ash Creek, 2012). The comments were provided to the Port of Portland (Port) in a letter from the DEQ dated July 24, 2012. The DEQ comments are repeated (in italics) followed by the Port response.

# Risk Assessment

- 1. Current guidance documents should have been used.
  - The latest DEQ human health risk assessment guidance is from 2010, not 2003. Current guidance supports the use of Risk-Based Concentrations (RBCs) to calculate risk. The screening procedure for multiple chemicals has been simplified.
  - The most current table of RBCs is from 2011, not 2009. The current RBCss for occupational exposure to PCBs is 0.56 mg/kg, not 0.98 mg/kg. For TPH-diesel, RBCss Occupational is 14,000 mg/kg, and the RBCss Construction Worker.
  - The current version of EPA's ProUCL is 4.1, not 4.00.04. UCL values developed using the older program appear to be reasonable.

Based on our review the use of current RBC values will not change the list of chemicals of potential concern (COPCs) or the risk calculations for arsenic and carcinogenic PAHs based on current RBCs. Therefore, no changes are requested for this document, but future risk assessments submitted to DEQ should be based on current guidance.

Response. Noted.

2. Section 4.1.3 Soil Exposure Point Calculations, Page 13, second bullet. The report does not state how the 90UCL was chosen from the ProUCL output. It appears that the recommendations provided by ProUCL for the 95UCL were used for selecting the 90UCL. This approach is acceptable to DEQ, however the method for deciding how the 90UCL was selected should have been explained in the report. Future risk assessments submitted to DEQ should include an explanation.

Response. Noted.

3. Section 4.1.3 Soil Exposure Point Calculations, Page 13, third bullet. EPA does not recommend using ½ detection limit for non-detect values. For datasets that include non-detected values, the mean should be calculated using a method such as Kaplan-Meir. The results are likely not substantially different that the mean value reported in Table 2, therefore no changes are requested to this document. Future risk assessments submitted to DEQ should follow EPA's recommendation in dealing with non-detected values.

**Response.** Noted. For clarification, the 90UCL values were calculated using Kaplan-Meir. The ½-detection limit method was used only for calculating mean values.

4. Section 4.3.1 Non-Carcinogenic Effect. The report states that there are no non-carcinogens identified as COPCs. DEQ notes that all chemicals, including those evaluated as carcinogens, have non-cancer effects. However, there are no reference doses available for carcinogenic PAHs, so non-cancer effects cannot be quantitatively evaluated. For arsenic, the lowest acceptable concentrations will be based on cancer effects. No changes to the report are requested.

Response. Noted.

5. **Figure 5 Human Health Conceptual Site Model.** Direct contact with soil should have explicitly shown that this includes incidental ingestion, dermal contact, and inhalation. For groundwater, these exposure routs are explicitly shown. However, the test on page 11 is clear that the different exposure routs included with "direct contact" were considered, therefore no changes to the report are requested.

Response. Noted.

6. **Table A-4. URS Soil Analytical Results.** The table does not present the method reporting limits, parameters, or screening level value included in the volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs), or polychlorinated biphenyls (PCBs). DEQ request that the tables be revised so that the adequacy of the method reporting limits can be compared to appropriate screening level values or that an evaluation of the adequacy of the method reporting limits be presented.

**Response.** A table presenting the range of method reporting limits (MRLs) for VOCs, SVOCs, and PCBs along with the current RBCss screening criteria is attached. Except for two PAHs, method reporting limits were less than the RBCs. PAHs were also analyzed by SIM methods with lower reporting limits so results are acceptable.

- 7. **Table A-8. URS Grab Groundwater Analytical Results.** The table does not present the method reporting limits, parameters, or screening level value included in the volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs), or polychlorinated biphenyls (PCBs). DEQ request that the tables be revised so that the adequacy of the method reporting limits can be compared to appropriate screening level values or that an evaluation of the adequacy of the method reporting limits be presented.
  - **Response.** A table presenting the range of method reporting limits (MRLs) for VOCs, SVOCs, and PCBs along with the JSCS screening criteria is attached. MRLs are generally consistent with industry standards (PAHs were also analyzed by SIM methods).
- 8. **Appendix C.** The first sets of ProUCL results for 0-15 feet do not appear to represent all the available data. We spot-checked evaluations of arsenic and benzo(a)pyrene, and were able to confirm the results for the subsequent 0-3 feet and 0-15 feet sets of UCL output sheets. Input values should have been explicitly presented (showing sample locations) so that the calculations could have been more easily confirmed.

Response. Noted.

9. **Appendix C.** The datasets included in the appendix show reporting limits for non-detected values. EPA and DEQ risk assessment guidance is to use estimated (J) values for detected concentrations that are below reporting limits. ProUCL is designed to incorporate non-detect values in estimating UCLs. It is inappropriate to substitute reporting limits for detection limits. Future submittals should report method detection limits and estimated values.

Response. Noted.

### Feasibility Study

10. **Section 5.1.1 Remedial Action Objectives.** The remedial action objectives (RAOs) are not clearly stated. RAOs should identify receptor(s) and pathway(s) that need to be addressed, e.g. 1) prevent site workers from direct contact of contaminated soil with concentrations exceeding RBCss Occupational or Construction Workers for benzo(a)pyrene 2) prevent site workers from direct contact of contaminated soil with concentrations exceeding RBCss Occupational Workers for arsenic.

Response. Noted.

11. **Section 5.4.2** Cap (Detailed Analysis of Remedial Action Alternatives). The capping alternative should identify development and implementation of a DEQ-approved Operations and Maintenance Plan (OMP) for the site to monitor and maintain the implemented remedy, in addition to the soil management plan (SMP). An OMP is somewhat covered in the SMP but should be identified in the report title (and procedures for implementation better-described). DEQ will provide comments on the SMP separately.

**Response.** This will be addressed in response to the DEQ comments on the SMP.

#### Source Control Evaluation

12. **Section 6.1.2 Chemicals of Interest.** DEQ requests that a figure showing the locations of sediment samples presented in Table D-1 through D6 be submitted.

**Response.** The noted sample locations are shown on Map 2.2-1k from LWG Draft Final RI report dated August 29, 2011. A copy of the Map is attached.

- 13. DEQ agrees with the findings and conclusions presented in the source control evaluation.
  - Groundwater is the only potential pathway associated with OU-4 (stormwater is a relevant pathway but will be evaluated by the current property operator).
  - The groundwater pathway does not appear to pose a recontamination risk to river sediments based on current soil, groundwater, and in-river sediment data.
  - Implementation of source control measures is not recommended for OU-4 at this time.

Response. Noted.

# **Next Steps**

Based on our review of the report, DEQ anticipates the following steps:

1. Port submit updated tables with reporting limits (or evaluation of the reporting limits) and a figure with sediment locations as requested above.

Response. Noted.

2. DEQ will propose implementation of source control measures are not needed at this time and document this in the form of a Draft Source Control Decision.

Response. Noted.

3. DEQ will submit the Draft Source Control Decision to EPA and the Portland Harbor Government team for review.

Response. Noted.

4. DEQ/Port will address EPA and Portland Harbor Government team review comments as needed.

Response. Noted.

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Please call me at (503) 415-6325 if you have any questions.

Sincerely,

**Dwight Leisle** 

Environmental Project Manager

c: Kristine Koch, EPA Rich Muza, EPA

Suzanne Barthelmess, Port

Richard Vincent, Port David Ashton, Port

Michael Pickering, Ash Creek Associates Mark Lewis, Formation Environmental

LWP File

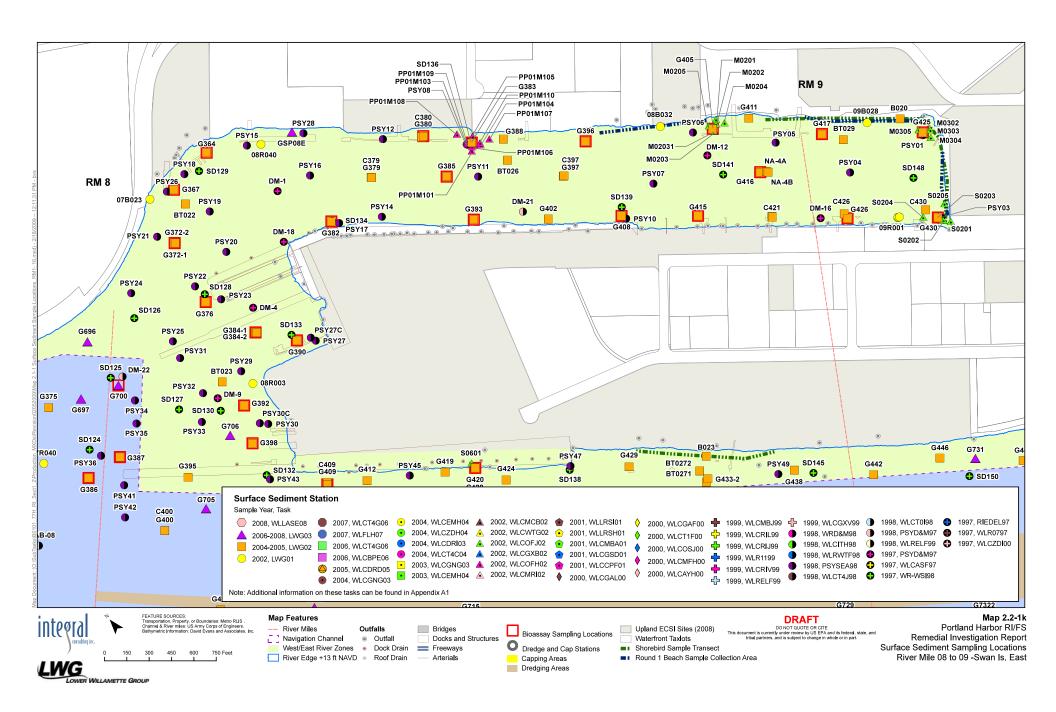


Table 1 - Analytical Detection Limits Summary - PCBs/Butyl Tins/SVOCs/VOCs URS OU4, Swan Island Upland Facility Portland, Oregon

Polychlorinated Biphenyls (PCBs) by EPA Method 8082	Portland, Oregon Sample Number	Water	JSCS	Soil	RBC
PCB-1016					INDO
PCB-1271	r oryomormacou Espironyio (r o			ug/kg	ug/kg
PCB-1232	PCB-1016	-	-		
PCB-1242		-	0.034		
PGB-1246					
PCB-1256					
PGB-1260					
Upl.					560
1.2.4-Trichlorobenzene   55   49   330 to 3300   300   19,000,1	Semi-volatile Organic Compour				ua/ka
1.2-Dichlorobenzene 1.2-Diphenylhydrazine 1.3-Dichlorobenzene 1.4-Dichlorobenzene 1.4-Dichlorobenzene 1.4-Dichlorobenzene 1.4-Dichlorobenzene 1.5-Dichlorobenzene 1.4-Dichlorobenzene 1.4-Dichlorobenzene 1.4-Dichlorobenzene 1.4-Dichloropheno 1.5-Dichloropheno 1.5-Di	1.2.4-Trichlorobenzene				
1,2-Diphenylyhdrazine					19,000,000
1.4-Dichlorobenzene	1,2-Diphenylhydrazine	<10		<330 to <3300	
2.4.5-Trichlorophenol	1,3-Dichlorobenzene	<5		<330 to <3300	
2.4.6-Trichlorophenol					63,000
2.4-Dinitrophenol	-				-
2.4-Dimitrophenol         <5					
2.4-Dinitrotoluene         <25					-
2.4-Dinitrotoluene         <5					
2.6-Dinitrotouene 2-5					
2-Chloronaphthalene 2-Chlorophenol 5-S 30 30 330 to 3300 3300 330 to 3300 34-Chlorophenyl ether 4-Chlorophenyl phenyl ether 4-Chlorophenyl phenyl ether 4-Chlorophenyl phenyl ether 4-Chlorophenyl phenyl ether 4-Nitrophenol 4-Chlorophenyl phenyl ether 4-Nitrophenol 4-Chlorophenyl ether 4-Nitrophenol 4-Nitrophenol 4-See to 20 330 to 3300 300 300 300 300 300 300 300 300 30	-				240,000
2-Chloropheno    <5					
2-Methylnaphthalene				<330 to <3300	-
2-Nitropheno		<5	0.2	<330 to <3300	
3,3 - Dichlorobenzidine	2-Nitroaniline	<25	110	<800 to <8000	-
3-Nitroaniline   <25			150	<330 to <3300	
4,6-Dinitro-2-methylphenol   4-Ehromophenyl phenyl ether   4-Chloros-methylphenol   4-Chloros-methylphenol   4-Chloropaniline					
4-Bromophenyl phenyl ether 4-Chloro-3-methylphenol 5			3.2		-
4-Chloro-3-methylphenol         <5			-		
4-Chloropaniline					
4-Chlorophenyl phenyl ether 4-Nitrophenol					
A-Nitroaniline					
A-Nitropheno					
Acenaphthylene Aniline <5	4-Nitrophenol			<800 to <8000	
Aniline Anthracene		<5		<330 to <3300	19,000,000
Anthracene   Senz(a)anthracene   Senzo(a)pyrene   Senzo(a)pyrene   Senzo(a)pyrene   Senzo(b)fluoranthene   Senzo(g,h,i)perylene   Senzo(g,h,i,i)perylene   Senzo(g,h,i,i,i)perylene   Senzo(g,h,i,i)perylene   Senzo(g,h,i,i)perylene   Senzo(g,h,i,i)perylene   Senzo					
Benz(a)anthracene   Senzo(a)pyrene   Senzo(a)pyrene   Senzo(b)fluoranthene   Senzo(b)fluo					
Benzo(a)pyrene   Senzo(b)fluoranthene   Sen					93,000,000
Benzo(b)fluoranthene					
Benzo(g,h,i)perylene					
Benzo(k)fluoranthene					
Benzoic acid Benzyl alcohol   <5   8.6   <330 to <3300					27,000
Bis(2-chloroethoxy)methane   Sis(2-chloroethyl)ether   Sis(2-chlorosopropyl)ether   Sis(2-chlorosopro	` '				
Bis(2-chloroethyl)ether   Sis(2-chloroisopropyl)ether   Sis(2-ch	Benzyl alcohol	<5	8.6	<330 to <3300	-
Bis(2-chloroisopropyl)ether   Sis(2-ethylhexyl)phthalate   Sis(2-ethylhe				<330 to <3300	
Bis(2-ethylhexyl)phthalate         < 5	,				
Butyl benzyl phthalate					450,000
Carbazole					150,000
Chrysene         < 5         0.018         <330 to <3300         250,00           Dibenz(a,h)anthracene         < 5					
Dibenz(a,h)anthracene         < 5         0.018         <330 to <3300         270           Dibenzofuran         < 5					250,000
Dibenzofuran         < 5	,				
Dimethyl phthalate         < 5		< 5			-
Di-n-butyl phthalate         < 5	Diethyl phthalate	< 5		<330 to <3300	
Di-n-octyl phthalate         < 5				<330 to <3300	
Fluoranthene Fluorene					
Fluorene					
Hexachlorobenzene         < 5					8,900,000
Hexachlorobutadiene         < 5					12,000,000
Hexachlorocyclopentadiene         < 5					1,200
Hexachloroethane					
Indeno(1,2,3-cd)pyrene					90,000
Isophorone					
Naphthalene       < 5					
Nitrobenzene         < 5	Naphthalene				23,000
N-Nitrosodiphenylamine       < 5	Nitrobenzene				
Pentachlorophenol         < 25         0.56         < 800 to < 8000         3,900           Phenanthrene         < 5					-
Phenanthrene       < 5					-
Phenol       < 5       2,560       <330 to <3300          Pyrene       < 5					3,900
Pyrene       < 5					-
Pyridine < 5 <330 to <3300 2-Methylphenol < 5 <330 to <3300					6 700 000
2-Methylphenol < 5 <330 to <3300			0.2		ο,/υυ,υυυ
3 & 4-Methylphenol < 5 <330 to <3300	3 & 4-Methylphenol				

Sample Number	Water	JSCS	Soil	RBC
Volatile Organic Compounds (V	OCs) by EPA Method			
	ug/L	ug/L	mg/kg	mg/kg
Acetone	<20.0	1,500	<0.944 to <1.52	
Benzene	<0.250	1.2	<0.0118 to <0.0190	34.0
Bromobenzene	< 0.500		<0.0236 to <0.0381	
Bromochloromethane	< 0.500		<0.0236 to <0.0381	
Bromodichloromethane	< 0.500	1.1	<0.0236 to <0.0381	15.0
Bromoform	<1.00	8.5	<0.0472 to <0.0761	240.0
Bromomethane	<5.00	8.7	<0.472 to <0.761	330
2-Butanone (MEK)	<10.0		<0.472 to <0.761	
n-Butylbenzene	<1.00		<0.0236 to <0.0381	
sec-Butylbenzene	<1.00		<0.0236 to <0.0381	
tert-Butylbenzene	<0.500		<0.0236 to <0.0381	
Carbon tetrachloride	<0.500	0.51	<0.0236 to <0.0381	31
Chlorobenzene	<0.500	50	<0.0236 to <0.0381	4,300
Chloroethane	<2.00	23	<0.472 to <0.761	>100,000
Chloroform	<2.00	0.17	<0.236 to <0.381	25
Chloromethane	<5.00	2.1	<0.236 to <0.381	25,000
2-Chlorotoluene	<0.500		<0.0236 to <0.0381	
4-Chlorotoluene	<0.500		<0.0236 to <0.0381	
1,2-Dibromo-3-chloropropane	<2.00	_	<0.0944 to <0.152	
Dibromochloromethane	<0.500	0.79	<0.0472 to <0.0761	16
1,2-Dibromoethane (EDB)	<0.500		<0.0236 to <0.0381	0.68
Dibromomethane	< 0.500	61	<0.0236 to <0.0381	
1,2-Dichlorobenzene	<0.500	49	<0.0236 to <0.0381	19,000
1,3-Dichlorobenzene	<0.500	14	<0.0236 to <0.0381	13,000
1,4-Dichlorobenzene	<0.500	2.8	<0.0236 to <0.0381	63
Dichlorodifluoromethane	<1.00	390	<0.0472 to <0.0761	03
1,1-Dichloroethane	<0.500	J90 	<0.0236 to <0.0381	250
	<0.500		<0.0236 to <0.0381	15
1,2-Dichloroethane (EDC)	<0.500		<0.0236 to <0.0381	
1,1-Dichloroethene cis-1,2-Dichloroethene	<0.500		<0.0236 to <0.0381	12,000 620
			<0.0236 to <0.0381	
trans-1,2-Dichloroethene	<0.500 <0.500		<0.0236 to <0.0381	4,500
1,2-Dichloropropane 1,3-Dichloropropane	<0.500		<0.0236 to <0.0381	
2,2-Dichloropropane	<0.500		<0.0236 to <0.0381	
1,1-Dichloropropene	<0.500		<0.0236 to <0.0381	
cis-1,3-Dichloropropene	<0.500	0.055	<0.0472 to <0.0761	
trans-1,3-Dichloropropene	<0.500	0.055	<0.0472 to <0.0761	
Ethylbenzene	<0.500	7.3	<0.0236 to <0.0381	140
Hexachlorobutadiene	<2.00	0.86	<0.0230 to <0.0361 <0.0944 to <0.152	140
2-Hexanone	<10.0	0.00	<0.472 to <0.761	
Isopropylbenzene	<0.500	660	<0.0236 to <0.0381	24,000
4-Isopropyltoluene	<0.500		<0.0236 to <0.0381	24,000
4-Methyl-2-pentanone (MiBK)	<0.500 <10.0		<0.472 to <0.761	
Methyl tert-butyl ether (MTBE)	<0.500	37	<0.472 to <0.761	
Methylene chloride	<5.00	8.9	<0.236 to <0.381	1,000
Netrylerie Chloride Napthalene	<5.00 <5.00	0.9	<0.236 to <0.381	23
n-Propylbenzene	<0.500		<0.0236 to <0.0381	
Styrene	<0.500	100	<0.0236 to <0.0381	51,000
1,1,1,2-Tetrachloroethane	<0.500	100	<0.0472 to <0.0761	51,000
1,1,2,2-Tetrachloroethane	<0.500 <0.500	_	<0.0472 to <0.0761 <0.0236 to <0.0381	
Tetrachloroethene (PCE)	<0.500	0.12	<0.0236 to <0.0381	940
Tetrachioroethene (PCE) Toluene	<0.500 <1.00	9.8	<0.0236 to <0.0361 <0.0944 to <0.152	24,000
		3.0	<0.0944 to <0.152	24,000
1,2,3-Trichlorobenzene 1,2,4-Trichlorobenzene	<2.00 <2.00	8.2	<0.0944 to <0.152	
1,2,4-Trichloroethane	<0.500	0.2	<0.0944 to <0.152 <0.0472 to <0.0761	430,000
1,1,2-Trichloroethane	<0.500 <0.500	_	<0.0472 to <0.0761 <0.0236 to <0.0381	430,000 25
Trichloroethylene (TCE)		0.17	<0.0236 to <0.0381	25 46
	<0.500			
Trichlorofluoromethane	<1.00	1,300	<0.236 to <0.381	63,000
1,2,3,-Trichloropropane	<1.00		<0.0472 to <0.0761	2.000
1,2,4-Trimethylbenzene	<1.00	-	<0.0472 to <0.0761	2,000
1,3,5-Trimethylbenzene	<1.00	0.045	<0.0472 to <0.0761	3,100
Vinyl Chloride	<0.500	0.015	<0.0236 to <0.0381	4
m,p-Xylene o-Xylene	<1.00 <0.500	1.8	<0.0472 to <0.0761	19,000
u-∧yierie	<0.500	13	<0.0236 to <0.0381	19,000

- 1. JSCS -- DEQ/EPA, 2005. Portland Harbor Joint Source Control Strategy Final (Table 3-1 Updated
- July 16, 2007). December 2005.

  2. RBC = DEQ risk-based concentration for lower of construction worker or occupational direct contact (June 2012 update). Value for PCBs is for total. 3. -- = Not applicable or not available.
- 4. μg/L = micrograms per liter.
- 5.  $\mu$ g/kg = micrograms per kilogram.